

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: ssptansc1625

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

NEWS	1		Web Page for STN Seminar Schedule - N. America
NEWS	2	JAN 02	STN pricing information for 2008 now available
NEWS	3	JAN 16	CAS patent coverage enhanced to include exemplified prophetic substances
NEWS	4	JAN 28	USPATFULL, USPAT2, and USPATOLD enhanced with new custom IPC display formats
NEWS	5	JAN 28	MARPAT searching enhanced
NEWS	6	JAN 28	USGENE now provides USPTO sequence data within 3 days of publication
NEWS	7	JAN 28	TOXCENTER enhanced with reloaded MEDLINE segment
NEWS	8	JAN 28	MEDLINE and LMEDLINE reloaded with enhancements
NEWS	9	FEB 08	STN Express, Version 8.3, now available
NEWS	10	FEB 20	PCI now available as a replacement to DPCI
NEWS	11	FEB 25	IFIREF reloaded with enhancements
NEWS	12	FEB 25	IMSPRODUCT reloaded with enhancements
NEWS	13	FEB 29	WPINDEX/WPIDS/WPIX enhanced with ECLA and current U.S. National Patent Classification
NEWS	14	MAR 31	IFICDB, IFIPAT, and IFIUDB enhanced with new custom IPC display formats
NEWS	15	MAR 31	CAS REGISTRY enhanced with additional experimental spectra
NEWS	16	MAR 31	CA/CAplus and CASREACT patent number format for U.S. applications updated
NEWS	17	MAR 31	LPCI now available as a replacement to LDPCI
NEWS	18	MAR 31	EMBASE, EMBAL, and LEMBASE reloaded with enhancements
NEWS	19	APR 04	STN AnaVist, Version 1, to be discontinued
NEWS	20	APR 15	WPIDS, WPINDEX, and WPIX enhanced with new predefined hit display formats
NEWS	21	APR 28	EMBASE Controlled Term thesaurus enhanced
NEWS	22	APR 28	TMSRESEARCH reloaded with enhancements

NEWS EXPRESS FEBRUARY 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 20 FEBRUARY 2008

NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS LOGIN Welcome Banner and News Items
NEWS IPC8 For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific

research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

FILE 'HOME' ENTERED AT 15:39:46 ON 04 MAY 2008

FILE 'REGISTRY' ENTERED AT 15:39:56 ON 04 MAY 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 2 MAY 2008 HIGHEST RN 1019057-98-7
DICTIONARY FILE UPDATES: 2 MAY 2008 HIGHEST RN 1019057-98-7

New CAS Information Use Policies. enter HELP USAGETERMS for details.

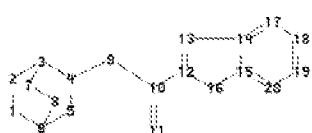
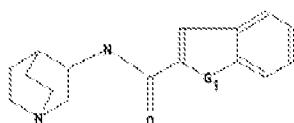
TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stnqen/stndoc/properties.html>

=>
Uploading C:\Program Files\Stnexp\Queries\10522611.str



```
chain nodes :
9 10 11
ring nodes :
1 2 3 4 5 6 7 8 12 13 14 15 16 17 18 19 20
chain bonds :
4-9 9-10 10-11 10-12
ring bonds :
```

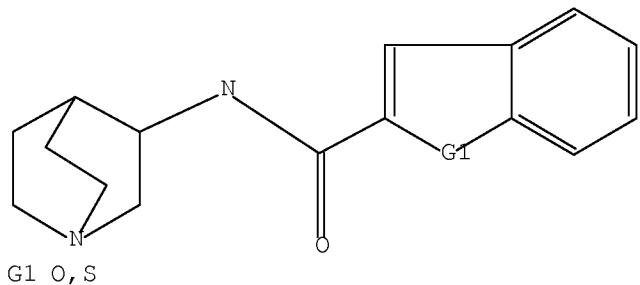
1-2 1-6 2-3 3-4 3-7 4-5 5-6 6-8 7-8 12-13 12-16 13-14 14-17 14-15 15-
16 15-20 17-18 18-19 19-20
exact/norm bonds :
1-2 1-6 2-3 3-4 3-7 4-5 4-9 5-6 6-8 7-8 9-10 10-11 10-12 12-13 12-16
13-14 15-16
normalized bonds :
14-17 14-15 15-20 17-18 18-19 19-20

G1:O, S

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:CLASS
11:CLASS 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
20:Atom

L1 STRUCTURE UPLOADED

=> d 11
L1 HAS NO ANSWERS
L1 STR



Structure attributes must be viewed using STN Express query preparation.

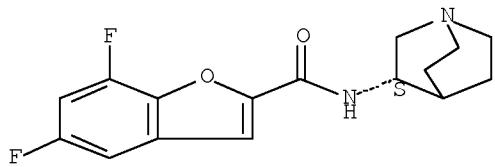
=> s sss sam 11
SAMPLE SEARCH INITIATED 15:40:18 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 42 TO ITERATE
100.0% PROCESSED 42 ITERATIONS 27 ANSWERS
SEARCH TIME: 00.00.01
FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 452 TO 1228
PROJECTED ANSWERS: 229 TO 851

L2 27 SEA SSS SAM L1

=> d scan

L2 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN 2-Benzofurancarboxamide, N-(3S)-1-azabicyclo[2.2.2]oct-3-yl-5,7-difluoro-
MF C16 H16 F2 N2 O2

Absolute stereochemistry.

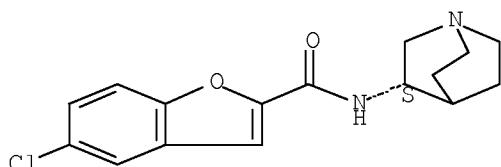


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):4

L2 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN 2-Benzofurancarboxamide, N-(3S)-1-azabicyclo[2.2.2]oct-3-yl-5-chloro-,
monohydrochloride (9CI)
MF C16 H17 Cl N2 O2 . Cl H

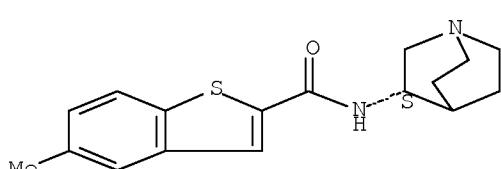
Absolute stereochemistry.



● HCl

L2 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Benzo[b]thiophene-2-carboxamide, N-(3S)-1-azabicyclo[2.2.2]oct-3-yl-5-
methyl-
MF C17 H20 N2 O S
CI COM

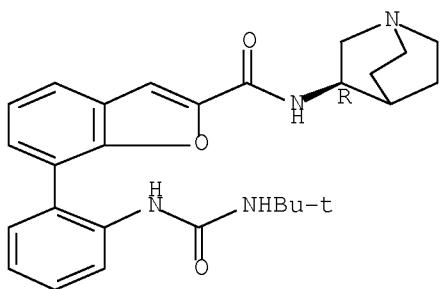
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN 2-Benzofurancarboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-7-[2-[[[(1,1-dimethylethyl)amino]carbonyl]amino]phenyl]-
MF C27 H32 N4 O3
CI COM

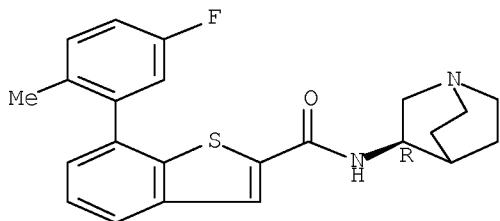
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-7-(5-fluoro-2-methylphenyl)-
MF C23 H23 F N2 O S
CI COM

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> s sss full 11

FULL SEARCH INITIATED 15:41:07 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 831 TO ITERATE

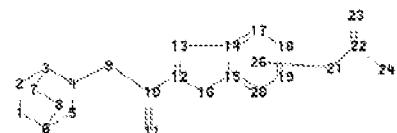
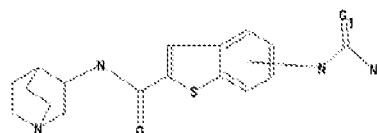
100.0% PROCESSED 831 ITERATIONS
SEARCH TIME: 00.00.01

647 ANSWERS

L3 647 SEA SSS FUL L1

=> save 13 fles10522611/a
ANSWER SET L3 HAS BEEN SAVED AS 'FLES10522611/A'

=>
Uploading C:\Program Files\Stnexp\Queries\10522611A.str



chain nodes :

9 10 11 23

ring nodes :

1 2 3 4 5 6 7 8 12 13 14 15 16 17 18 19 20

ring/chain nodes :

21 22 24

chain bonds :

4-9 9-10 10-11 10-12 21-22 22-23 22-24

ring bonds :

1-2 1-6 2-3 3-4 3-7 4-5 5-6 6-8 7-8 12-13 12-16 13-14 14-17 14-15 15-16 15-20 17-18 18-19 19-20

exact/norm bonds :

1-2 1-6 2-3 3-4 3-7 4-5 4-9 5-6 6-8 7-8 9-10 10-11 12-13 12-16 13-14 15-16 21-22 22-23 22-24

exact bonds :

10-12

normalized bonds :

14-17 14-15 15-20 17-18 18-19 19-20

G1:O,S

Match level :

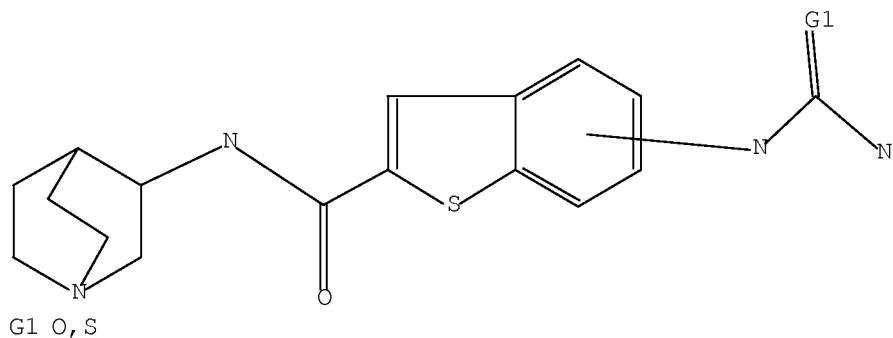
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:CLASS
11:CLASS 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
20:Atom 21:CLASS 22:CLASS 23:CLASS 24:CLASS 26:Atom

L4 STRUCTURE UPLOADED

=> D L4

L4 HAS NO ANSWERS

L4 STR



Structure attributes must be viewed using STN Express query preparation.

=> S SSS SAM SUBSET=L3
 ENTER LOGIC EXPRESSION, QUERY NAME, OR (END):END
 SEARCH ENDED BY USER

=> S SSS SAM SUBSET=L3 L4
 SAMPLE SUBSET SEARCH INITIATED 15:47:10 FILE 'REGISTRY'
 SAMPLE SUBSET SCREEN SEARCH COMPLETED - 2 TO ITERATE

100.0% PROCESSED 2 ITERATIONS 2 ANSWERS
 SEARCH TIME: 00.00.01

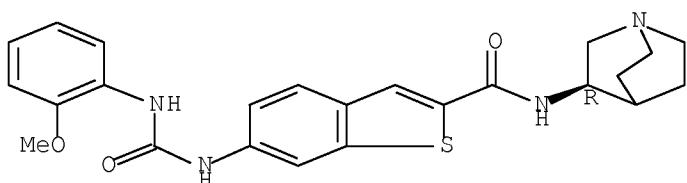
PROJECTIONS (WITHIN SPECIFIED SUBSET):	ONLINE	**COMPLETE**
PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET):	2 TO	124
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET):	2 TO	124

L5 2 SEA SUB=L3 SSS SAM L4

=> D SCAN

L5 2 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-
 [[[2-methoxyphenyl]amino]carbonyl]amino]-, monohydrochloride (9CI)
 MF C24 H26 N4 O3 S . Cl H

Absolute stereochemistry.



● HCl

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> S SSS FULL SUBSET=L3 L4
FULL SUBSET SEARCH INITIATED 15:47:28 FILE 'REGISTRY'
FULL SUBSET SCREEN SEARCH COMPLETED - 44 TO ITERATE

100.0% PROCESSED 44 ITERATIONS 42 ANSWERS
SEARCH TIME: 00.00.01

L6 42 SEA SUB=L3 SSS FUL L4

=> FIL CAPLU
COST IN U.S. DOLLARS SINCE FILE TOTAL
FULL ESTIMATED COST ENTRY SESSION
225.98 226.19

FILE 'CAPLUS' ENTERED AT 15:47:38 ON 04 MAY 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 4 May 2008 VOL 148 ISS 19
FILE LAST UPDATED: 2 May 2008 (20080502/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/infopolicy.html>

=> S L6
L7 1 L6

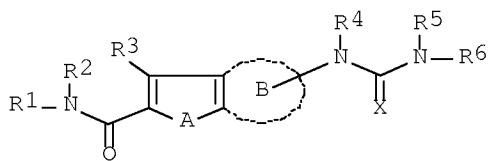
=> D IBIB ABS HITSTR

L7 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2004:117115 CAPLUS Full-text
DOCUMENT NUMBER: 140:181328
TITLE: Preparation of benzothiophenes, benzofuranes and indoles for improvement of cognition, attention achievement, learning achievement and/or memory achievement
INVENTOR(S): Flessner, Timo; Boess, Frank-Gerhard; Erb, Christina; Hafner, Frank-Thorsten; Schnizler, Katrin; Lang, Dieter; Luithle, Joachim; Van Kampen, Marja; Van der Staay, Franz-Josef

PATENT ASSIGNEE(S): Bayer AG, Germany
 SOURCE: Ger. Offen., 35 pp.
 CODEN: GWXXBX
 DOCUMENT TYPE: Patent
 LANGUAGE: German
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
DE 10234424	A1	20040212	DE 2002-10234424	20020729
CA 2494352	A1	20040212	CA 2003-2494352	20030714
WO 2004013136	A1	20040212	WO 2003-EP7588	20030714
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
AU 2003254343	A1	20040223	AU 2003-254343	20030714
EP 1527073	A1	20050504	EP 2003-766162	20030714
EP 1527073	B1	20070822		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK				
JP 2005538111	T	20051215	JP 2004-525191	20030714
ES 2291694	T3	20080301	ES 2003-766162	20030714
US 20060106096	A1	20060518	US 2005-522611	20051215
PRIORITY APPLN. INFO.:			DE 2002-10234424	A 20020729
			WO 2003-EP7588	W 20030714

OTHER SOURCE(S): MARPAT 140:181328
 GI



AB Title compds. [I; R1 = 1-azabicyclo[2.2.2]oct-3-yl; R2 = H, alkyl; R3 = H, halo, amino, OH, C1-6 alkyl; R4 = H, (substituted) C1-6 alkyl; R5 = H, C1-6 alkyl; or R4R5 = (substituted) 5-6 membered heterocyclyl; R6 = H, alkyl, cycloalkyl, aryl, (substituted) 5-6 membered heterocyclyl, etc.; NR5R6 = (substituted) 3-8 membered heterocyclyl; A = O, N, S; X = O, S; B = (substituted) benzo or pyrido ring], were prepared. Thus, 6-amino-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-1-benzothiophene-2-carboxamide dihydrochloride (preparation given) in THF was reacted with Et3N and 2,6-difluorophenyl isocyanate for 18 h at room temperature to give 75.9% N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-6-[(2,6-difluorophenyl)aminocarbonyl]aminocarbonylamine-1-benzothiophene-2-carboxamide hydrochloride. The latter showed affinity to α 7-NACHR by inhibiting [3H]methyllycaconite binding in rat brain membranes with $K_i = 0.4$ nM.

IT 657401-24-6P 657401-25-7P 657401-26-8P
657401-27-9P 657401-28-0P 657401-29-1P
657401-30-4P 657401-31-5P 657401-32-6P
657401-33-7P 657401-34-8P 657401-36-0P
657401-37-1P 657401-38-2P 657401-39-3P
657401-40-6P 657401-41-7P 657401-42-8P
657401-43-9P 657401-44-0P 657401-45-1P

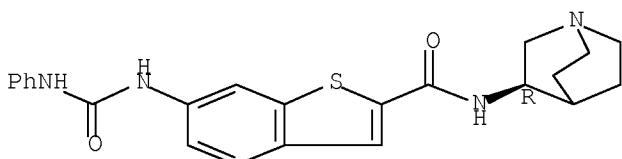
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of benzothiophenes, benzofuranes and indoles for improvement of cognition, attention achievement, learning achievement and/or memory achievement)

RN 657401-24-6 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[(phenylamino)carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

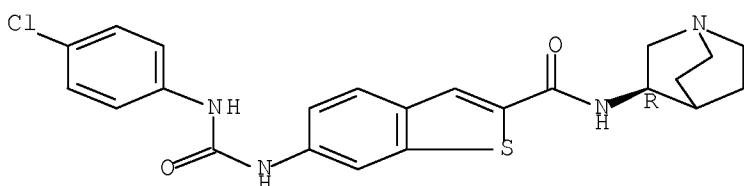


● HCl

RN 657401-25-7 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[(4-chlorophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

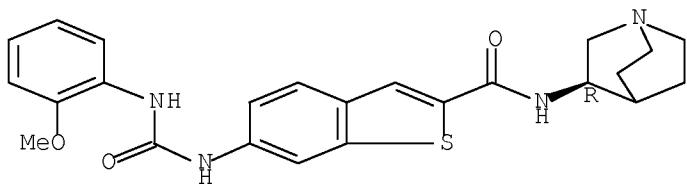


● HCl

RN 657401-26-8 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[(2-methoxyphenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

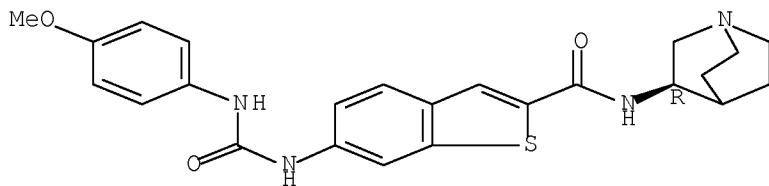


● HCl

RN 657401-27-9 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-
[[(4-methoxyphenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA
INDEX NAME)

Absolute stereochemistry.

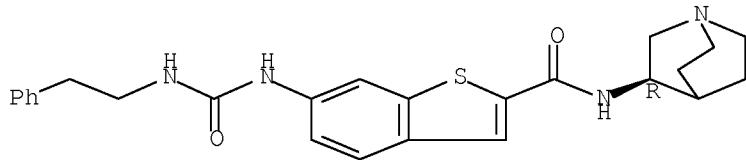


● HCl

RN 657401-28-0 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-
[[(2-phenylethyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA
INDEX NAME)

Absolute stereochemistry.

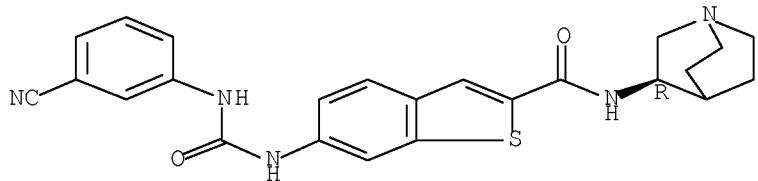


● HCl

RN 657401-29-1 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-
[[(3-cyanophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA
INDEX NAME)

Absolute stereochemistry.

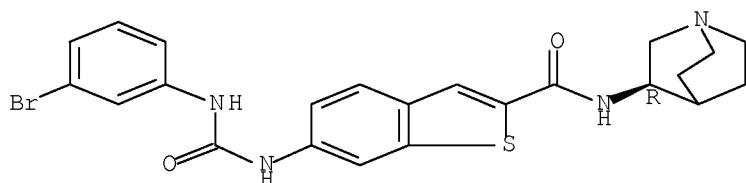


● HCl

RN 657401-30-4 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(3-bromophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

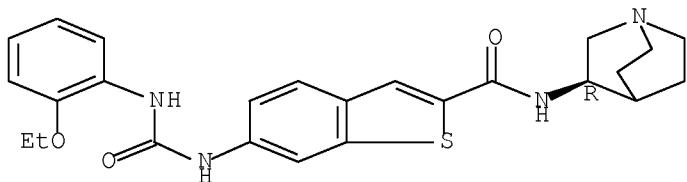


● HCl

RN 657401-31-5 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(2-ethoxyphenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

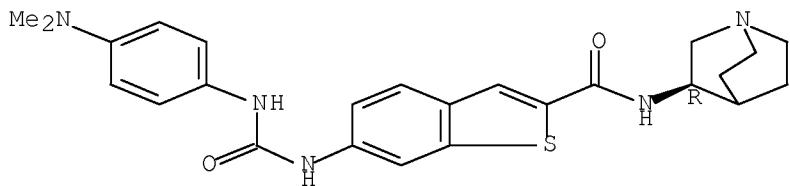


● HCl

RN 657401-32-6 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(4-(dimethylamino)phenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

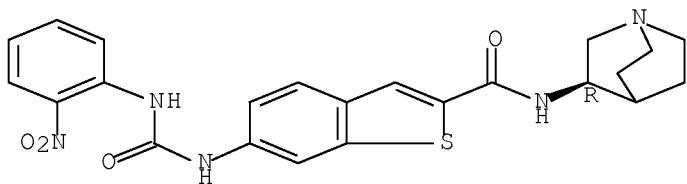


● HCl

RN 657401-33-7 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(2-nitrophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

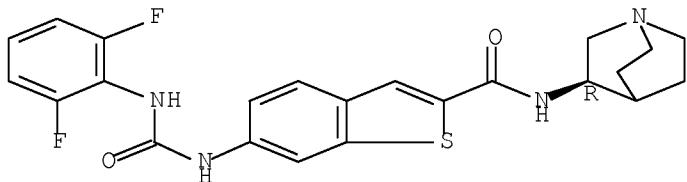


● HCl

RN 657401-34-8 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(2,6-difluorophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

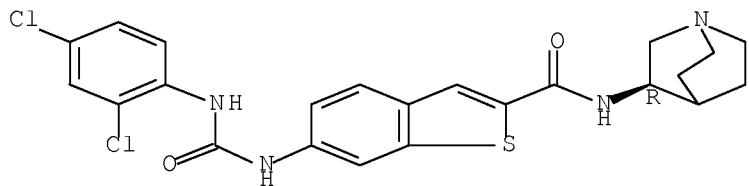


● HCl

RN 657401-36-0 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(2,4-dichlorophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

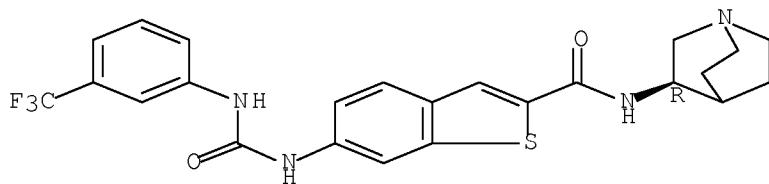


● HCl

RN 657401-37-1 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[3-(trifluoromethyl)phenyl]amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

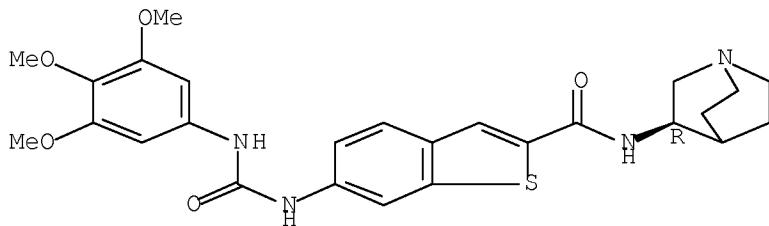


● HCl

RN 657401-38-2 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[3,4,5-trimethoxyphenyl]amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

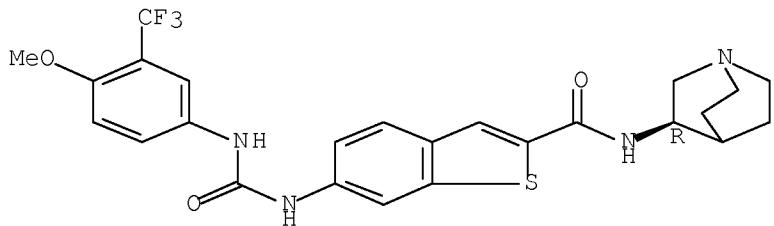


● HCl

RN 657401-39-3 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[4-methoxy-3-(trifluoromethyl)phenyl]amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

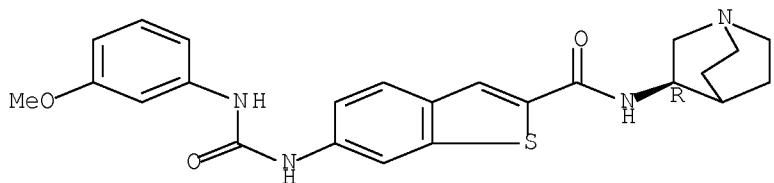


● HCl

RN 657401-40-6 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[(3-methoxyphenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

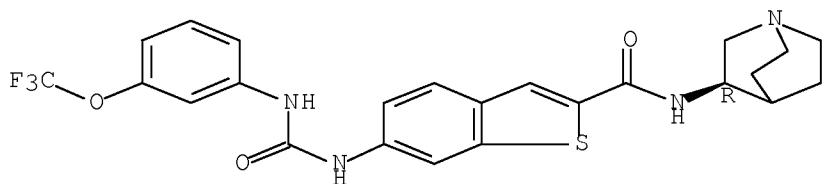


● HCl

RN 657401-41-7 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[(3-(trifluoromethoxy)phenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

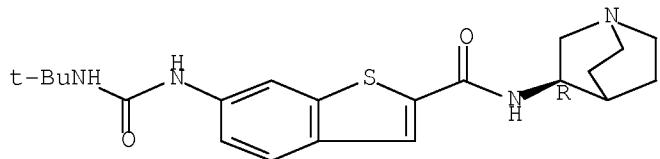


● HCl

RN 657401-42-8 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[(1,1-dimethylethyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

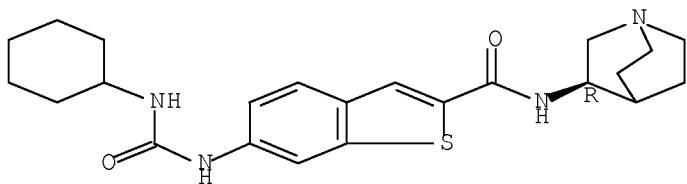


● HCl

RN 657401-43-9 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[(cyclohexylamino)carbonyl]amino-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

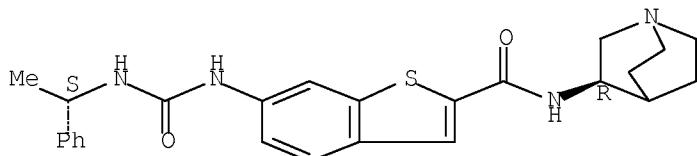


● HCl

RN 657401-44-0 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(1S)-1-phenylethyl]amino]carbonyl]amino-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

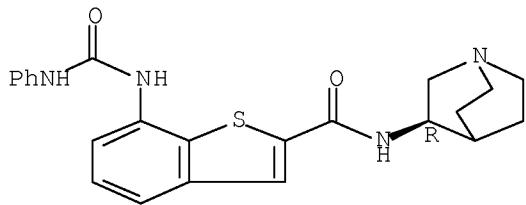


● HCl

RN 657401-45-1 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-7-[(phenylamino)carbonyl]amino-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.



● HCl

=> LOGOFF Y
COST IN U.S. DOLLARS
FULL ESTIMATED COST

	SINCE FILE ENTRY	TOTAL SESSION
	5.93	232.12

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)
CA SUBSCRIBER PRICE

	SINCE FILE ENTRY	TOTAL SESSION
	-0.80	-0.80

STN INTERNATIONAL LOGOFF AT 15:47:57 ON 04 MAY 2008